

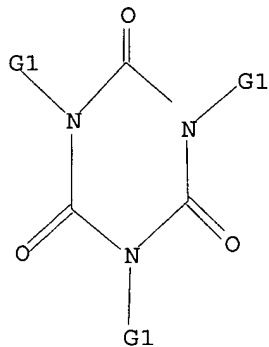
Uploading 09863463.str

L8 STRUCTURE UPLOADED

=> d l8

L8 HAS NO ANSWERS

L8 STR



G1 X, Ak, H

Structure attributes must be viewed using STN Express query preparation.

=> s l8

SAMPLE SEARCH INITIATED 16:08:03 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 71 TO ITERATE

100.0% PROCESSED 71 ITERATIONS

6 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**

BATCH **COMPLETE**

PROJECTED ITERATIONS: 915 TO 1925

PROJECTED ANSWERS: 6 TO 266

L9 6 SEA SSS SAM L8

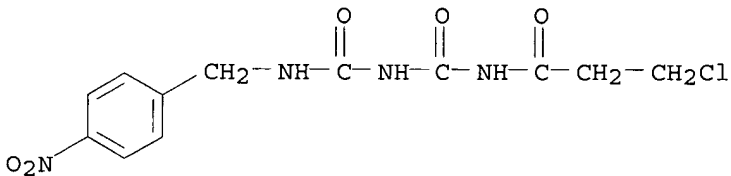
=> d scan

L9 6 ANSWERS REGISTRY COPYRIGHT 2003 ACS

IN Propanamide, 2-chloro-N-[[[(4-nitrophenyl)methyl]amino]carbonyl]amino]carbonyl]- (9CI)

MF C12 H13 Cl N4 O5

CI COM



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):end

=> s l8 full

FULL SEARCH INITIATED 16:08:23 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 1616 TO ITERATE

100.0% PROCESSED 1616 ITERATIONS 169 ANSWERS
SEARCH TIME: 00.00.01

L10 169 SEA SSS FUL L8

=> file caplus

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	148.55	305.85

FILE 'CAPLUS' ENTERED AT 16:08:31 ON 19 MAR 2003
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2003 AMERICAN CHEMICAL SOCIETY (ACS)

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FILE COVERS 1907 - 19 Mar 2003 VOL 138 ISS 12
FILE LAST UPDATED: 18 Mar 2003 (20030318/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s l8 and electrolyte

REGISTRY INITIATED

Substance data SEARCH and crossover from CAS REGISTRY in progress...
Use DISPLAY HITSTR (or FHITSTR) to directly view retrieved structures.

SAMPLE SEARCH INITIATED 16:08:39 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 71 TO ITERATE

100.0% PROCESSED 71 ITERATIONS 6 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS:	ONLINE	**COMPLETE**
	BATCH	**COMPLETE**
PROJECTED ITERATIONS:	915 TO	1925
PROJECTED ANSWERS:	6 TO	266

L11 6 SEA SSS SAM L8

L12 10 L11

209573 ELECTROLYTE
L13 0 L12 AND ELECTROLYTE

=> s l10 and electrolyte
69 L10

209573 ELECTROLYTE
L14 0 L10 AND ELECTROLYTE

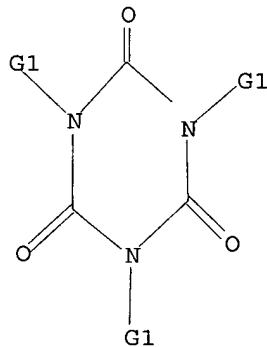
=> s l10 and battery
69 L10

97692 BATTERY
L15 0 L10 AND BATTERY

=>
Uploading 09863463.str

L1 STRUCTURE UPLOADED

=> d l1
L1 HAS NO ANSWERS
L1 STR



G1 X, Ak, H

Structure attributes must be viewed using STN Express query preparation.

=> s l1
SAMPLE SEARCH INITIATED 16:17:09 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 71 TO ITERATE

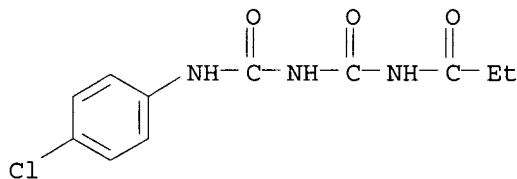
100.0% PROCESSED 71 ITERATIONS 6 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
 BATCH **COMPLETE**
PROJECTED ITERATIONS: 915 TO 1925
PROJECTED ANSWERS: 6 TO 266

L2 6 SEA SSS SAM L1

=> d scan 1-6
'1-6' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'

L2 6 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Propanamide, N-[[[(4-chlorophenyl)amino]carbonyl]amino]carbonyl] - (9CI)
MF C11 H12 Cl N3 O3



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

The following are valid formats:

Substance information can be displayed by requesting individual fields or predefined formats. The predefined substance formats are: (RN = CAS Registry Number)

REG - RN
SAM - Index Name, MF, and structure - no RN
FIDE - All substance data, except sequence data
IDE - FIDE, but only 50 names
SQIDE - IDE, plus sequence data
SQIDE3 - Same as SQIDE, but 3-letter amino acid codes are used
SQD - Protein sequence data, includes RN
SQD3 - Same as SQD, but 3-letter amino acid codes are used
SQN - Protein sequence name information, includes RN

CALC - Table of calculated properties
EPROP - Table of experimental properties
PROP - EPROP and CALC

Any CA File format may be combined with any substance format to obtain CA references citing the substance. The substance formats must be cited first. The CA File predefined formats are:

ABS -- Abstract
APPS -- Application and Priority Information
BIB -- CA Accession Number, plus Bibliographic Data
CAN -- CA Accession Number
CBIB -- CA Accession Number, plus Bibliographic Data (compressed)
IND -- Index Data
IPC -- International Patent Classification
PATS -- PI, SO
STD -- BIB, IPC, and NCL

IABS --ABS, indented, with text labels
IBIB -- BIB, indented, with text labels
ISTD -- STD format, indented

OBIB ----- AN, plus Bibliographic Data (original)
OIBIB ----- OBIB, indented with text labels

SBIB ----- BIB, no citations
SIBIB ----- IBIB, no citations

The ALL format gives FIDE BIB ABS IND RE, plus sequence data when it is available.

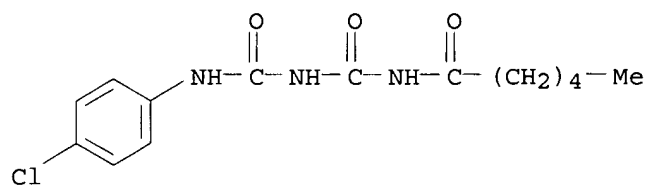
The MAX format is the same as ALL.

The IALL format is the same as ALL with BIB ABS and IND indented, with text labels.

For additional information, please consult the following help messages:

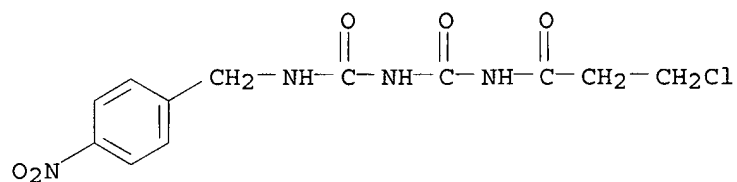
HELP DFIELDS -- To see a complete list of individual display fields.
HELP FORMATS -- To see detailed descriptions of the predefined formats.
HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):5

L2 6 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Hexanamide, N-[[[(4-chlorophenyl)amino]carbonyl]amino]carbonyl]- (9CI)
MF C14 H18 Cl N3 O3



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

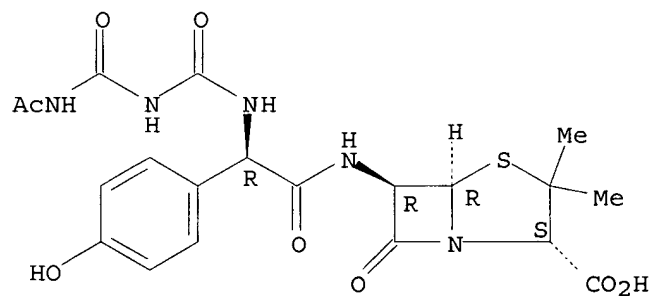
L2 6 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN . Propanamide, 2-chloro-N-[[[[(4-nitrophenyl)methyl]amino]carbonyl]amino]carbonyl]- (9CI)
 MF C12 H13 Cl N4 O5
 CI COM



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L2 6 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN 4-Thia-1-azabicyclo[3.2.0]heptane-2-carboxylic acid, 6-[[[[(4-hydroxyphenyl)acetyl]amino]carbonyl]amino] (4-hydroxyphenyl)acetyl]amino]-3,3-dimethyl-7-oxo-, monopotassium salt, [2S-[2.alpha.,5.alpha.,6.beta.(S*)]]- (9CI)
 MF C20 H23 N5 O8 S . K

Absolute stereochemistry.

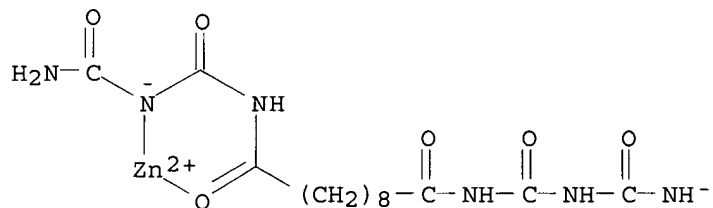


● K

$$\text{n-BuNH}-\overset{\text{O}}{\parallel}\text{C}-\text{NH}-\overset{\text{O}}{\parallel}\text{C}-\text{NH}-\overset{\text{O}}{\parallel}\text{C}-\text{Et}$$
$$\text{Me}-(\text{CH}_2)_4-\overset{\text{O}}{\parallel}\text{C}-\text{NH}-\overset{\text{O}}{\parallel}\text{C}-\text{NH}-\overset{\text{O}}{\parallel}\text{C}-\text{NH}-\overset{\text{O}}{\parallel}\text{C}-(\text{CH}_2)_4-\text{Me}$$
$$\text{H}_2\text{N}-\overset{\overset{\text{O}}{\parallel}}{\text{C}}-\overset{\overset{\text{O}}{\parallel}}{\text{N}}^--\text{Cu}^{2+}\text{O}=\text{C}-\text{NH}-\text{C}(\text{CH}_3)=\text{O}-(\text{CH}_2)_3-\overset{\overset{\text{O}}{\parallel}}{\text{C}}-\text{NH}-\overset{\overset{\text{O}}{\parallel}}{\text{C}}-\text{NH}-\overset{\overset{\text{O}}{\parallel}}{\text{C}}-\text{NH}-$$

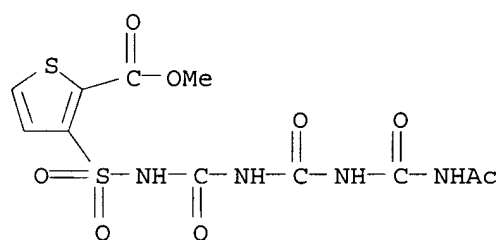
HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN Zinc, [3,14,16-trioxo-5-(oxo-.kappa.O)-2,4,15,17-tetraazaoctadecanediimidato(2-)-.kappa.N2]- (9CI)
 MF C14 H22 N6 O6 Zn
 CI CCS, COM



HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

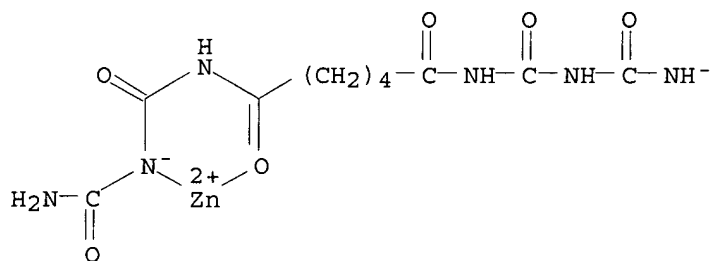
L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN 2-Thiophenecarboxylic acid, 3-[[[[[[[acetyl]amino]carbonyl]amino]carbonyl]amino]carbonyl]amino]sulfonyl]-, methyl ester (9CI)
 MF C11 H12 N4 O8 S2



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

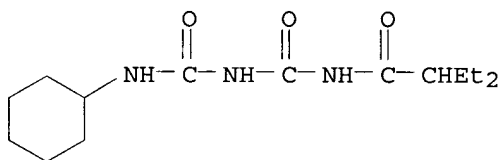
L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN Zinc, [3,5,10,12-tetraoxo-2,4,11,13-tetraazatetradecanediimidato(2-)-N2,O5]-, dihydrate (9CI)
 MF C10 H14 N6 O6 Zn . 2 H2 O
 CI CCS



● 2 H₂O

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

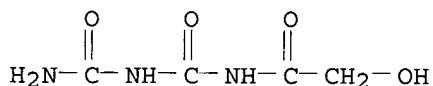
L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN Butanamide, N-[[[(cyclohexylamino)carbonyl]amino]carbonyl]-2-ethyl- (9CI)
 MF C14 H25 N3 O3



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

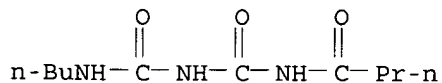
L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN Acetamide, N-[[[(aminocarbonyl)amino]carbonyl]-2-hydroxy- (9CI)
 MF C4 H7 N3 O4



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

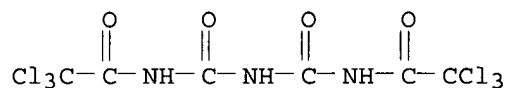
L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN Butanamide, N-[[[(butylamino)carbonyl]amino]carbonyl]- (9CI)
 MF C10 H19 N3 O3



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

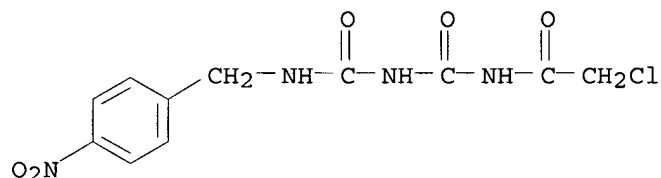
L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Acetamide, N,N'-(iminodicarbonyl)bis[2,2,2-trichloro- (9CI)
MF C6 H3 Cl6 N3 O4



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

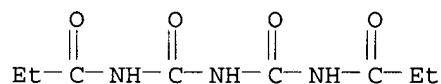
L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Acetamide, 2-chloro-N-[[[(4-nitrophenyl)methyl]amino]carbonyl]amino]carbonyl]- (9CI)
MF C11 H11 Cl N4 O5
CI COM



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Propanamide, N,N'-(iminodicarbonyl)bis- (9CI)
MF C8 H13 N3 O4
CI COM

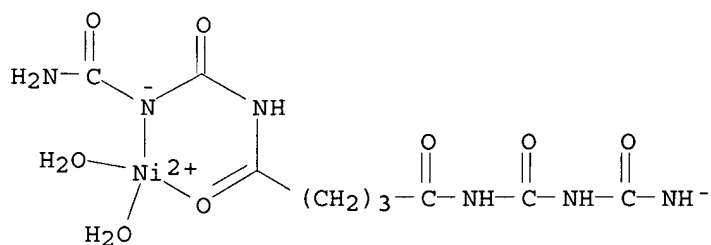


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

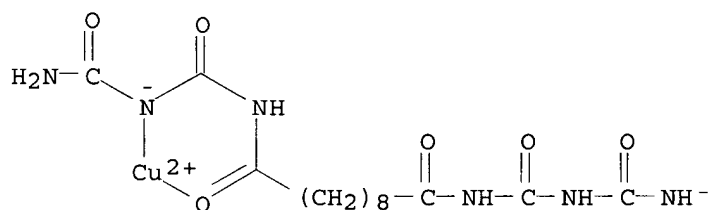
L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Nickel, diaqua[3,9,11-trioxo-5-(oxo-.kappa.O)-2,4,10,12-tetraazatridecanediamidato(2-)-.kappa.N2]- (9CI)

MF C9 H16 N6 Ni O8
CI CCS



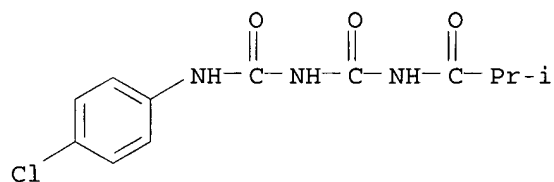
HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Copper, [3,14,16-trioxo-5-(oxo-.kappa.O)-2,4,15,17-tetraazaoctadecanediimidato(2-)-.kappa.N2]- (9CI)
MF C14 H22 Cu N6 O6
CI CCS



HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Propanamide, N-[[[(4-chlorophenyl)amino]carbonyl]amino]carbonyl]-2-methyl- (9CI)
MF C12 H14 Cl N3 O3

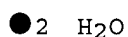
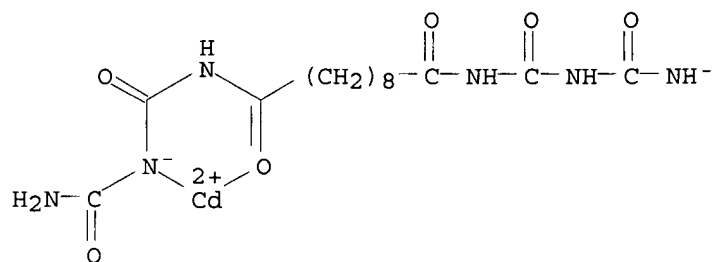


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

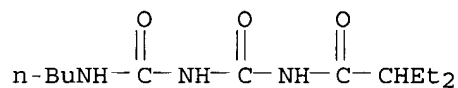
L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Cadmium, [3,5,14,16-tetraoxo-2,4,15,17-tetraazaoctadecanediimidato(2-)-N2,O5]-, dihydrate (9CI)

MF C14 H22 Cd N6 O6 . 2 H2 O
 CI CCS



HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

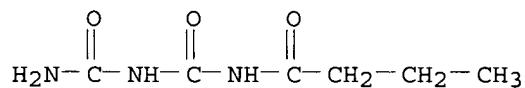
L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN Butanamide, N-[[[(butylamino)carbonyl]amino]carbonyl]-2-ethyl- (9CI)
 MF C12 H23 N3 O3



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

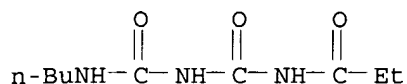
L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN Butanamide, N-[[[(aminocarbonyl)amino]carbonyl]- (9CI)
 MF C6 H11 N3 O3



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

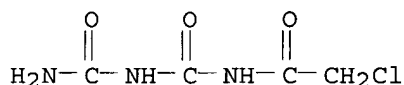
L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN Propanamide, N-[[[(butylamino)carbonyl]amino]carbonyl]- (9CI)
 MF C9 H17 N3 O3



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

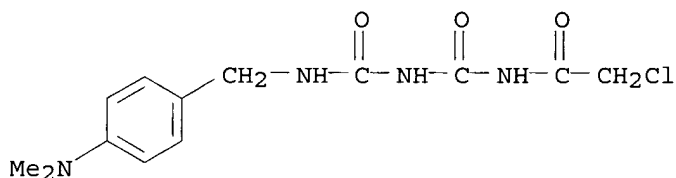
L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN Acetamide, N-[[[(aminocarbonyl)amino]carbonyl]-2-chloro- (9CI)
 MF C4 H6 Cl N3 O3



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

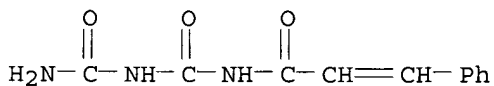
L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN Acetamide, 2-chloro-N-[[[[[4-(dimethylamino)phenyl]methyl]amino]carbonyl]amino]carbonyl]- (9CI)
 MF C13 H17 Cl N4 O3
 CI COM



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

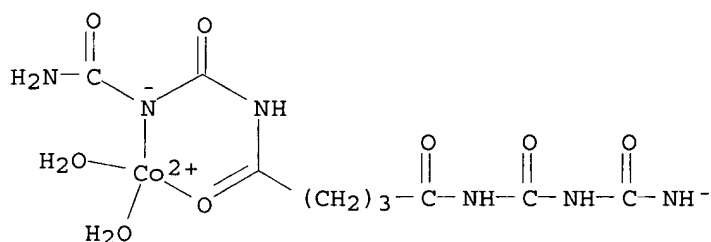
L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN 2-Propenamide, N-[[[(aminocarbonyl)amino]carbonyl]-3-phenyl- (9CI)
 MF C11 H11 N3 O3



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

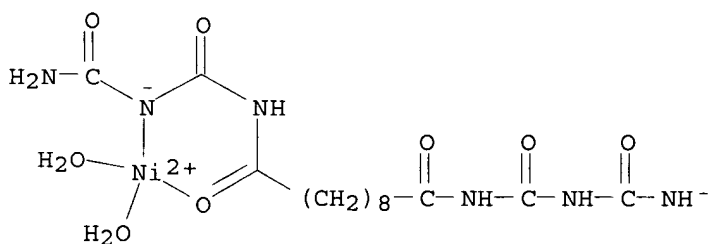
HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN Cobalt, diaqua[3,9,11-trioxo-5-(oxo-.kappa.O)-2,4,10,12-tetraazatridecanediamidato(2-)-.kappa.N2]-, monohydrate (9CI)
 MF C9 H16 Co N6 O8 . H2 O
 CI CCS



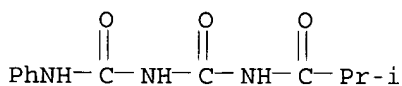
HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN Nickel, diaqua[3,14,16-trioxo-5-(oxo-.kappa.O)-2,4,15,17-tetrazaoctadecanediamidato(2-)-.kappa.N2]- (9CI)
 MF C14 H26 N6 Ni O8
 CI CCS



HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

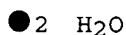
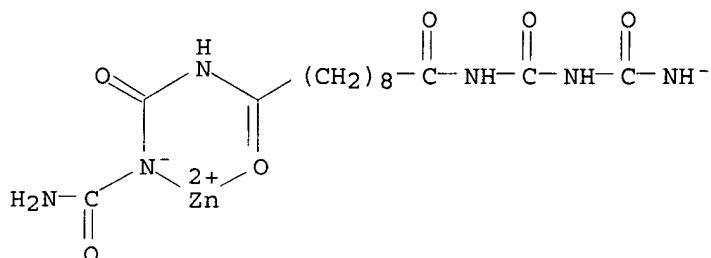
L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN Propanamide, 2-methyl-N-[[[(phenylamino)carbonyl]amino]carbonyl]- (9CI)
 MF C12 H15 N3 O3



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

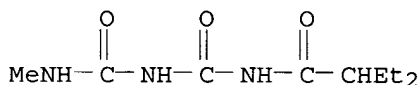
HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Zinc, [3,5,14,16-tetraoxo-2,4,15,17-tetraazaoctadecanediimidato(2-)-N2,O5]-
, dihydrate (9CI)
MF C14 H22 N6 O6 Zn . 2 H2 O
CI CCS



HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

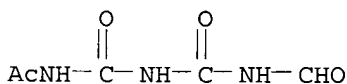
L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Butanamide, 2-ethyl-N-[[[(methylamino)carbonyl]amino]carbonyl]- (9CI)
MF C9 H17 N3 O3



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Biuret, 1-acetyl-5-formyl- (7CI)
MF C5 H7 N3 O4

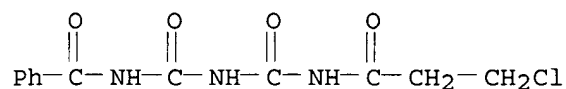


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Benzamide, N-[[[(3-chloro-1-oxopropyl)amino]carbonyl]amino]carbonyl]-
(9CI)

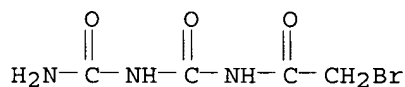
MF C12 H12 Cl N3 O4



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):
HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Acetamide, N-[[[(aminocarbonyl)amino]carbonyl]-2-bromo- (9CI)
MF C4 H6 Br N3 O3

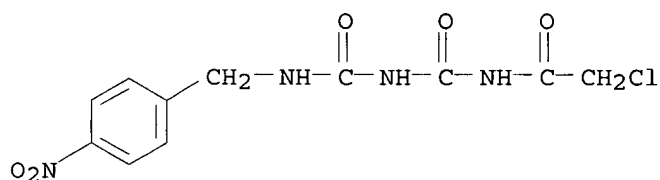


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

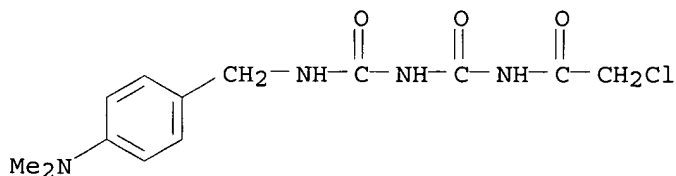
HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Acetamide, 2-chloro-N-[[[[[4-(dimethylamino)phenyl]methyl]amino]carbonyl]amino]carbonyl]-, compd. with 2-chloro-N-[[[[[4-nitrophenyl]methyl]amino]carbonyl]amino]carbonyl]acetamide (1:1) (9CI)
MF C13 H17 Cl N4 O3 . C11 H11 Cl N4 O5

CM 1

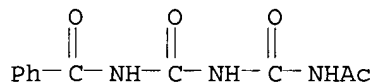


CM 2



HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

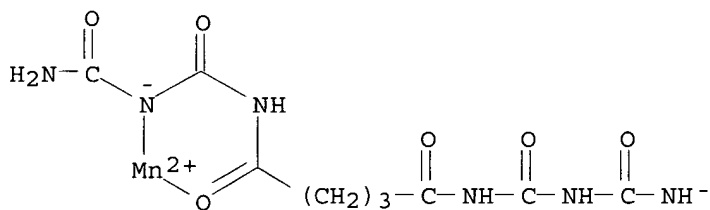
MF C11 H11 N3 O4



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

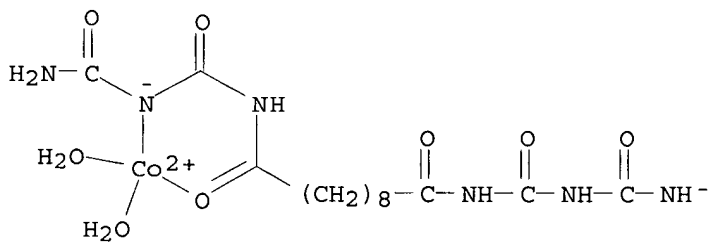
HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

CI CCS



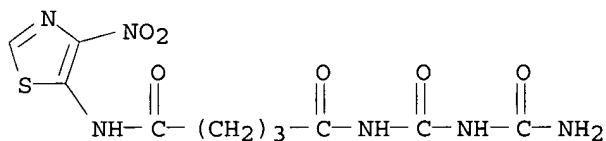
HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

CI CCS

 H_2O

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1) : 1

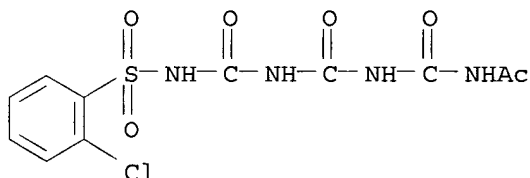
L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN Pentanediamide, N-[[[(aminocarbonyl)amino]carbonyl]-N'-(4-nitro-5-thiazolyl)- (9CI)
 MF C10 H12 N6 O6 S



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

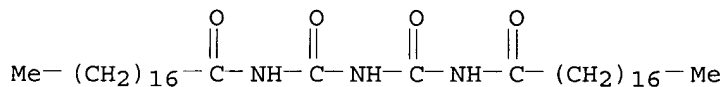
L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN Acetamide, N-[[[[[[[(2-chlorophenyl)sulfonyl]amino]carbonyl]amino]carbonyl]amino]carbonyl]- (9CI)
 MF C11 H11 Cl N4 O6 S



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

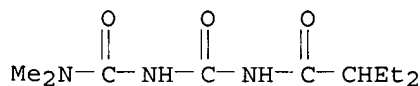
L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN Biuret, 1,5-distearoyl- (7CI)
 MF C38 H73 N3 O4

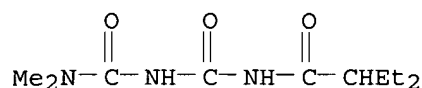


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN Butanamide, N-[[[(dimethylamino)carbonyl]amino]carbonyl]-2-ethyl- (9CI)
 MF C10 H19 N3 O3

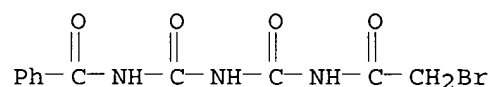




PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

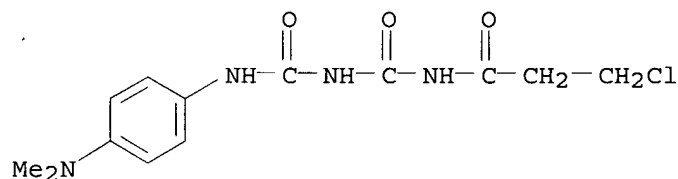
L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN Benzamide, N-[[[(bromoacetyl)amino]carbonyl]amino]carbonyl] - (9CI)
 MF C11 H10 Br N3 O4



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN Propanamide, 3-chloro-N-[[[4-(dimethylamino)phenyl]amino]carbonyl]amino]carbonyl] - (9CI)
 MF C13 H17 Cl N4 O3



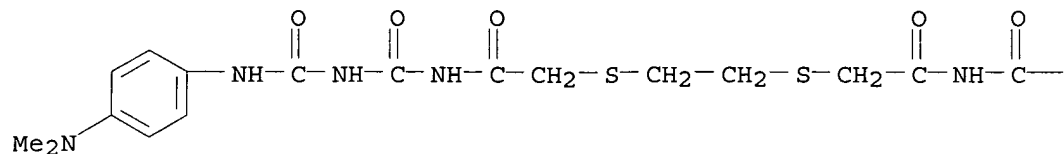
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

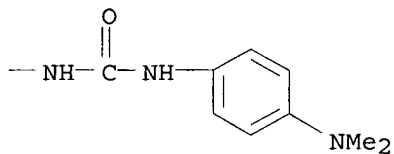
L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN 7,10-Dithia-2,4,13,15-tetraazahexadecanedi-2,13-diamide, N,N'-bis[4-(dimethylamino)phenyl]-3,5,12,14-tetraoxo-, compd. with N,N'-bis[(4-nitrophenyl)acetyl]-3,12-dioxo-6,9-dithia-2,13-diazatetradecanedi-2,13-diamide (1:1) (9CI)
 MF C26 H34 N8 O6 S2 . C26 H28 N6 O10 S2

CM 1

PAGE 1-A

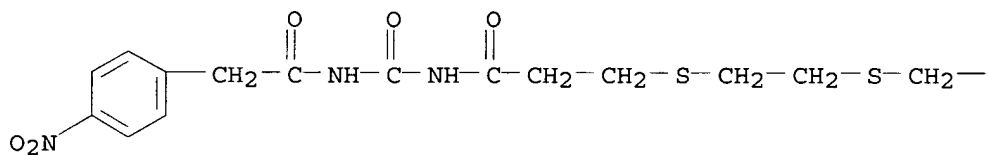


PAGE 1-B

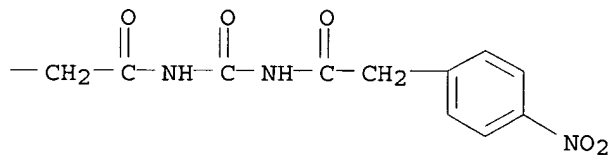


CM 2

PAGE 1-A

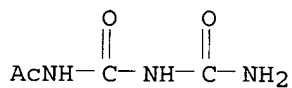


PAGE 1-B



HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

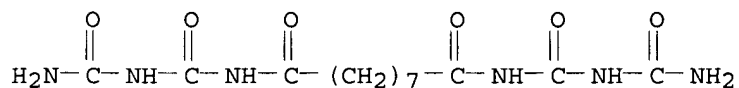
L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Acetamide, N-[[(aminocarbonyl) amino] carbonyl] - (9CI)
MF C4 H7 N3 O3



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

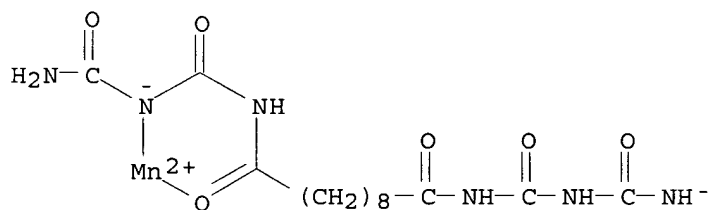
L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN 2,4,14,16-Tetraazaheptadecanediamide, 3,5,13,15-tetraoxo- (9CI)
MF C13 H22 N6 O6



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

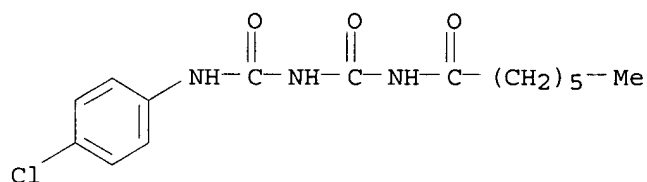
HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN Manganese, [3,14,16-trioxo-5-(oxo-.kappa.O)-2,4,15,17-tetraazaoctadecanediimidato(2-)-.kappa.N2]- (9CI)
 MF C14 H22 Mn N6 O6
 CI CCS



HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

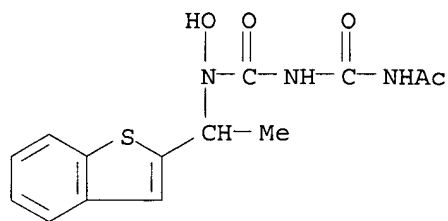
L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN Heptanamide, N-[[[(4-chlorophenyl)amino]carbonyl]amino]carbonyl]- (9CI)
 MF C15 H20 Cl N3 O3



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

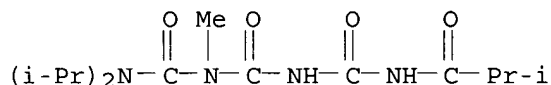
L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN Acetamide, N-[[[(1-benzo[b]thien-2-ylethyl)hydroxyamino]carbonyl]amino]carbonyl]- (9CI)
 MF C14 H15 N3 O4 S



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

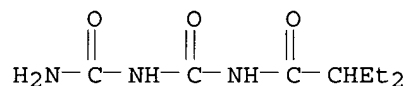
L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN Propanamide, N-[[[[(bis(1-methylethyl)amino)carbonyl]methylamino]carbonyl
]amino]carbonyl]-2-methyl- (9CI)
 MF C14 H26 N4 O4



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

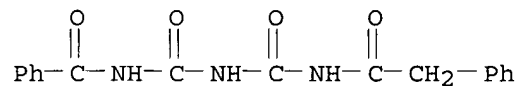
L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN Butanamide, N-[(aminocarbonyl)amino]carbonyl]-2-ethyl- (9CI)
 MF C8 H15 N3 O3



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN Benzeneacetamide, N-[[[(benzoylamino)carbonyl]amino]carbonyl]- (9CI)
 MF C17 H15 N3 O4

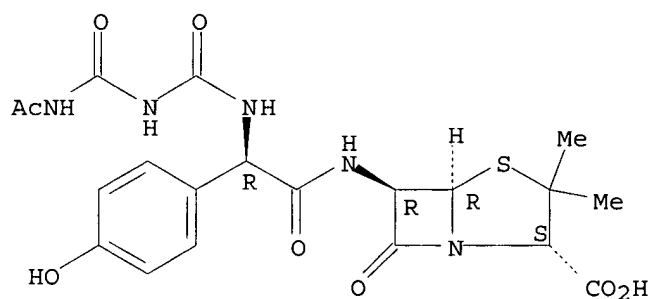


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN 4-Thia-1-azabicyclo[3.2.0]heptane-2-carboxylic acid, 6-
 [[[[[(acetylamino)carbonyl]amino]carbonyl]amino] (4-
 hydroxyphenyl)acetyl]amino]-3,3-dimethyl-7-oxo-, monopotassium salt,
 [2S-[2.alpha.,5.alpha.,6.beta.(S*)]]- (9CI)
 MF C20 H23 N5 O8 S . K

Absolute stereochemistry.



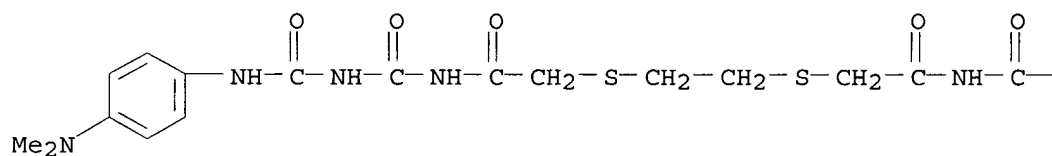
● K

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

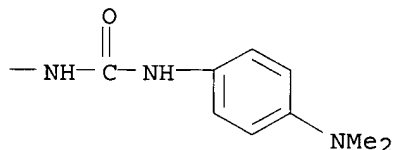
L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN 7,10-Dithia-2,4,13,15-tetraazahexadecanediamide, N,N'-bis[4-(dimethylamino)phenyl]-3,5,12,14-tetraoxo-, compd. with
 N,N'-bis[(4-nitrophenyl)acetyl]-3,10-dioxo-5,8-dithia-2,11-diazadodecanediamide (1:1) (9CI)
 MF C26 H34 N8 O6 S2 . C24 H24 N6 O10 S2

CM 1

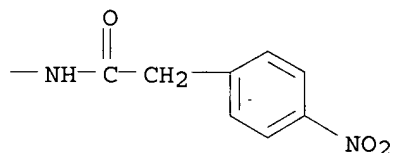
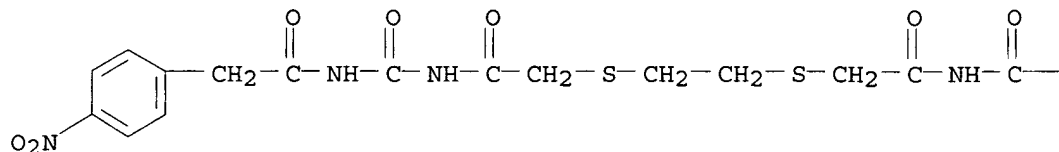
PAGE 1-A



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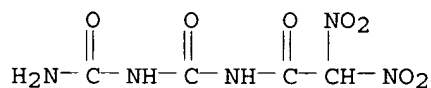


CM 2



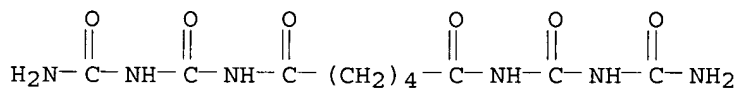
HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN Acetamide, N-[[(aminocarbonyl) amino] carbonyl]-2,2-dinitro-, monoammonium salt (9CI)
 MF C4 H5 N5 O7 . H3 N



HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN 2,4,11,13-Tetraazatetradecanediamide, 3,5,10,12-tetraoxo- (9CI)
 MF C10 H16 N6 O6

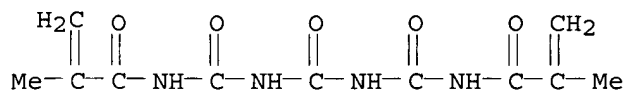


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

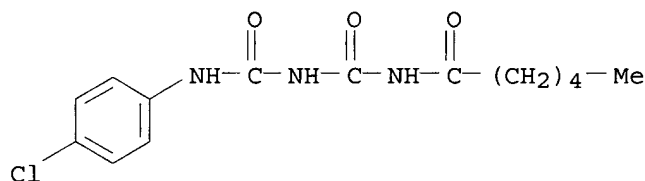
L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN 2-Propenamide, N,N'-[carbonylbis(iminocarbonyl)]bis[2-methyl-, homopolymer (9CI)
 MF (C11 H14 N4 O5)x
 CI PMS

CM 1



HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Hexanamide, N-[[[(4-chlorophenyl)amino]carbonyl]amino]carbonyl] - (9CI)
MF C14 H18 Cl N3 O3

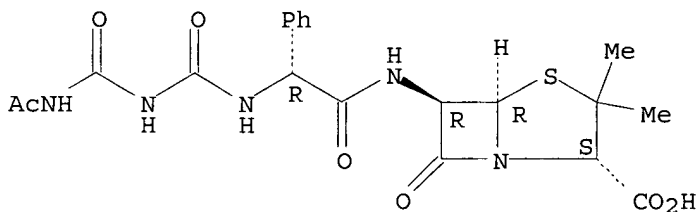


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN 4-Thia-1-azabicyclo[3.2.0]heptane-2-carboxylic acid, 6-
[[[[[(acetylamino)carbonyl]amino]carbonyl]amino]phenylacetyl]amino]-3,3-
dimethyl-7-oxo-, [2S-[2.alpha.,5.alpha.,6.beta.(S*)]] - (9CI)
MF C20 H23 N5 O7 S
CI COM

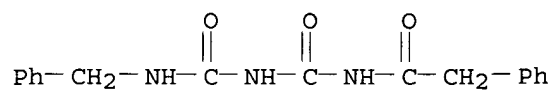
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

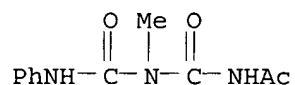
L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Benzeneacetamide, N-[[[(phenylmethyl)amino]carbonyl]amino]carbonyl] -
(9CI)
MF C17 H17 N3 O3



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L3 169 ANSWERS REGISTRY COPYRIGHT 2003 ACS
 IN Acetamide, N-[[methyl[(phenylamino)carbonyl]amino]carbonyl]- (9CI)
 MF C11 H13 N3 O3



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT